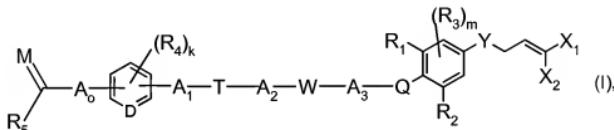


AMENDMENTS TO THE CLAIMS

Kindly amend claims 1 – 4 and cancel claims 8 – 9 without prejudice to the subject matter involved.
This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Currently amended) A compound of formula



wherein

A₀, A₁ and A₂ are each independently of the others a bond or a C₁-C₆alkylene bridge which is unsubstituted or substituted by from one to six identical or different substituents selected from halogen and C₃-C₈cycloalkyl;

A₃ is a C₁-C₆alkylene bridge which is unsubstituted or substituted by from one to six identical or different substituents selected from halogen and C₃-C₈cycloalkyl;

Y is O, NR₁₁, S, SO or SO₂;

M is O or NOR₆,

X₁ and X₂ are each independently of the other fluorine, chlorine or bromine;

R₁, R₂ and R₃ are each independently of the others H, halogen, OH, SH, CN, nitro, C₁-C₆alkyl, C₁-C₆haloalkyl, C₁-C₆alkylcarbonyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₂-C₆alkenyloxy, C₂-C₆haloalkenyloxy, C₂-C₆alkynyoxy, -S(=O)-C₁-C₆alkyl, -S(=O)₂-C₁-C₆alkyl, C₁-C₆alkoxycarbonyl or C₃-C₆haloalkynyoxy; the substituents R₃ being independent of one another when m is 2;

Q is O, NR₁₁, S, SO or SO₂;

W is O, NR₁₁, S, SO, SO₂, -C(=O)-O-, -O-C(=O)-, -C(=O)-NR₁₁- or -NR₁₁-C(=O)-;

T is a bond, O, NR₁₁, S, SO, SO₂, -C(=O)-O-, -O-C(=O)-, -C(=O)-NR₁₁- or -NR₁₁-C(=O)-;

D is CH or N;

R_4 is H, halogen, OH, SH, CN, nitro, C_1 - C_6 alkyl, C_1 - C_6 haloalkyl, C_1 - C_6 alkylcarbonyl, C_2 - C_6 alkenyl, C_2 - C_6 haloalkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, C_1 - C_6 haloalkoxy, C_2 - C_6 alkenylxylo, C_2 - C_6 alkynylxylo, $-S(=O)-C_1-C_6$ alkyl, $-S(=O)_2-C_1-C_6$ alkyl, C_1 - C_6 alkoxycarbonyl, C_3 - C_6 haloalkynylxylo, NH_2 , $NH(C_1-C_6$ alkyl) or $N(C_1-C_6$ alkyl) $_2$ wherein the two alkyl groups are independent of one another; the substituents R_4 being independent of one another when k is greater than 1;

R_5 is C_1-C_12 alkoxy- C_1-C_12 alkyl or heterocyclyl; C_1-C_12 alkyl substituted by from one to five substituents selected from the group consisting of N_3 , NO_2 , OH, C_3-C_8 cycloalkyl, C_3-C_8 cycloalkoxy, C_1-C_6 alkoxy, C_1-C_6 haloalkoxy, C_2-C_6 alkenylxylo, C_2-C_6 haloalkenylxylo, C_3-C_6 alkynylxylo, C_3-C_6 haloalkynyl, C_3-C_6 haloalkynylxylo, C_3-C_6 cycloalkyl- C_1-C_6 alkoxy, C_4-C_6 alkylcarbonyl, C_1-C_6 alkoxy- C_1-C_6 alkoxy, $P(=O)(OC_1-C_6$ alkyl) $_2$, $S(O)_q$ - R_{13} , NH_2 , $NH(C_1-C_6$ alkyl), $N(C_1-C_6$ alkyl) $_2$, wherein the two alkyl groups are independent of one another, $N(R_7)_2$, wherein the two R_7 s are independent of one another and $NR_{14}S(O)_qR_{15}$;

C_3-C_6 cycloalkyl substituted by from one to five identical or different substituents selected from the group consisting of C_1-C_6 alkyl, halogen, CN, NO_2 , OH, C_1-C_6 alkoxy, C_1-C_6 haloalkoxy, NH_2 , $NH(C_1-C_6$ alkyl) and $N(C_1-C_6$ alkyl) $_2$, wherein the two alkyl groups are independent of one another, $N(R_7)_2$, wherein the two R_7 s are independent of one another;

$-C(=O)-O-R_6$; $-C(=O)-R_6$; $-C(=O)-NH-R_6$; $-C(=N-O-R_6)R_{16}$; $-C(=N-NH-R_6)R_{16}$; C_2-C_6 alkenyl; C_2-C_6 alkynyl; heterocyclyl; or

$-NR_{14}S(O)_qR_{15}$

wherein the alkenyl and alkynyl radicals are unsubstituted or, depending upon the possibilities of substitution, substituted by from one to five identical or different substituents selected from the group consisting of halogen, N_3 , CN, NO_2 , OH, C_3-C_8 cycloalkyl, C_1-C_6 alkoxy, C_1-C_6 haloalkoxy, C_2-C_6 alkenylxylo, C_2-C_6 haloalkenylxylo, C_3-C_6 alkynylxylo, C_3-C_6 haloalkynylxylo, C_3-C_6 cycloalkyl- C_1-C_6 alkoxy, C_1-C_6 alkylcarbonyl, C_1-C_6 haloalkylcarbonyl, C_1-C_6 alkoxycarbonyl, C_1-C_6 alkylcarbonyl- C_1-C_6 alkyl, C_1-C_6 alkoxycarbonyl- C_1-C_6 alkyl, C_3-C_6 haloalkynyl, C_1-C_6 alkoxy- C_1-C_6 alkyl, C_1-C_6 haloalkoxy- C_1-C_6 alkyl, C_2-C_6 alkenylxylo- C_1-C_6 alkyl, C_2-C_6 haloalkenylxylo- C_1-C_6 alkyl, C_1-C_6 alkynyl, C_1-C_6 alkynylxylo- C_1-C_6 alkyl, $P(=O)(OC_1-C_6$ alkyl) $_2$, $S(O)_q$ - R_{13} , NH_2 , $NH(C_1-C_6$ alkyl) and $N(C_1-C_6$ alkyl) $_2$, wherein the two alkyl groups are independent of one another;

and wherein the heterocyclyl radical mentioned under R_5 are unsubstituted or, depending upon the possibilities of substitution, substituted by from one to five substituents selected from

halogen, CN, NO₂, OH, SH, C₁-C₆alkyl, C₁-C₆haloalkyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₃-C₆alkynyl, C₃-C₆cycloalkyl, C₃-C₆cycloalkyl-C₁-C₆alkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₂-C₆alkenylxylo, C₂-C₆haloalkenylxylo, C₃-C₆alkynylxylo, C₃-C₆haloalkynylxylo, C₃-C₆cycloalkyl-C₁-C₆alkoxy, C₁-C₆alkylcarbonyl, C₁-C₆haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylcarbonyl, C₁-C₆alkyl, C₁-C₆alkoxycarbonyl-C₁-C₆alkyl, C₁-C₆alkylthio, C₂-C₆alkenylthio, C₃-C₆alkynylthio, C₃-C₆cycloalkyl-C₁-C₆alkylthio, C₃-C₆haloalkynyl, C₂-C₆haloalkenylthio, C₁-C₆haloalkylthio, C₁-C₆alkoxy-C₁-C₆alkyl, C₁-C₆haloalkoxy-C₁-C₆alkyl, C₂-C₆alkenylxylo-C₁-C₆alkyl, C₂-C₆haloalkenylxylo-C₁-C₆alkyl, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ wherein the two alkyl groups are independent of one another, C₁-C₆alkylcarbonylamino, C₁-C₆haloalkylcarbonylamino, C₁-C₆alkoxycarbonylamino and C₁-C₆alkylaminocarbonylamino;

or, when A₆ is a C₂-C₆alkylene bridge, R₅ is C₂-C₆alkylene bonded to one of the carbon atoms of A₆;

or, when R₄ and a group C(=NOR₆)R₅ are in the ortho-position relative to one another, R₄ and R₅ together form a C₂-C₆alkylene bridge wherein one or two CH₂ groups each independently of the ether may be replaced by O, NR₁₂, S or SO, and wherein the CH₂ groups are unsubstituted or mono- or di- substituted by halogen, OH, SH, CN, nitro, C₁-C₆alkyl, C₁-C₆haloalkyl, C₁-C₆alkoxy or C₁-C₆haloalkoxy;

R₆ is H, C₁-C₁₂alkyl, C₃-C₆cycloalkyl, C₁-C₆alkylcarbonyl, C₂-C₆alkenyl, C₂-C₆alkynyl, aryl, heterocycl or benzyl, wherein the alkyl, cycloalkyl, alkenyl and alkynyl radicals are unsubstituted or, depending upon the possibilities of substitution, substituted by from one to five identical or different substituents selected from the group consisting of halogen, -N₃, CN, NO₂, OH, SH, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₂-C₆alkenylxylo, C₂-C₆haloalkenylxylo, C₃-C₆alkynylxylo, C₃-C₆cycloalkyl-C₁-C₆alkoxy, C₁-C₆alkylcarbonyl, C₁-C₆haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylcarbonyl-C₁-C₆alkyl, C₁-C₆alkylthio, C₂-C₆alkenylthio, C₃-C₆alkynylthio, C₃-C₆cycloalkyl-C₁-C₆alkylthio, C₃-C₆haloalkynyl, C₂-C₆haloalkenylthio, C₁-C₆haloalkylthio, C₁-C₆alkoxy-C₁-C₆alkyl, C₁-C₆haloalkoxy-C₁-C₆alkyl, C₂-C₆alkenylxylo-C₁-C₆alkyl, C₂-C₆haloalkenylxylo-C₁-C₆alkyl, C₃-C₆alkynylxylo-C₁-C₆alkyl, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ wherein the two alkyl groups are independent of one another, C₁-C₆alkylcarbonylamino, C₁-C₆haloalkylcarbonylamino, C₁-C₆alkoxycarbonylamino and C₁-C₆alkylaminocarbonylamino;

and the aryl, heterocycl and benzyl radicals are unsubstituted or, depending upon the possibilities of substitution, substituted by from one to five identical or different substituents selected

from the group consisting of halogen, CN, NO₂, OH, SH, C₁-C₆alkyl, C₁-C₆haloalkyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₃-C₆alkynyl, C₃-C₈cycloalkyl, C₃-C₈cycloalkyl-C₁-C₆alkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₂-C₆alkenylloxy, C₂-C₆haloalkenylloxy, C₃-C₆alkynylloxy, C₃-C₆haloalkynylloxy, C₃-C₈cycloalkyl-C₁-C₆alkoxy, C₁-C₆alkylcarbonyl, C₁-C₆haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylcarbonyl-C₁-C₆alkyl, C₁-C₆alkoxycarbonyl-C₁-C₆alkyl, C₁-C₆alkylthio, C₂-C₆alkenylthio, C₃-C₆alkynylthio, C₃-C₆cycloalkyl-C₁-C₆alkylthio, C₃-C₆haloalkynyl, C₂-C₆haloalkenylthio, C₁-C₆haloalkylthio, C₁-C₆alkoxy-C₁-C₆alkyl, C₁-C₆haloalkoxy-C₁-C₆alkyl, C₂-C₆alkenylloxy-C₁-C₆alkyl, C₂-C₆haloalkenylloxy-C₁-C₆alkyl, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ wherein the two alkyl groups are independent of one another, C₁-C₆alkylcarbonylamino, C₁-C₆haloalkylcarbonylamino, C₁-C₆alkoxycarbonylamino and C₁-C₆alkylaminocarbonylamino;

R₇ is H, C₁-C₆alkyl, C₁-C₈haloalkyl, C₁-C₆alkylcarbonyl, C₁-C₆haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₃-C₈cycloalkyl, C₃-C₈cycloalkylcarbonyl or formyl;

R₈ is H, C₁-C₁₂alkyl substituted by from one to five identical or different substituents selected from halogen, -N₃, CN, NO₂, OH, C₁-C₆alkoxy, C₁-C₆alkylthio, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ wherein the two alkyl groups are independent of one another and C₁-C₆alkylcarbonylamino; C₃-C₈cycloalkyl, C₁-C₆alkylcarbonyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₂-C₆alkynyl, C₂-C₆haloalkynyl, aryl, heterocycl or benzyl, wherein the aryl, heterocycl and benzyl radicals are unsubstituted or, depending upon the possibilities of substitution, substituted by from one to five substituents selected from the group consisting of halogen, CN, NO₂, OH, C₁-C₆haloalkyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₃-C₆alkynyl, C₃-C₈cycloalkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₂-C₆alkenylloxy, C₂-C₆haloalkenylloxy, C₃-C₆alkynylloxy, C₃-C₆haloalkynylloxy, C₁-C₆alkylcarbonyl, C₁-C₆haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylthio, C₂-C₆alkenylthio, C₃-C₆alkynylthio, C₁-C₆alkylthio, C₃-C₆haloalkynyl, C₁-C₆haloalkylthio, C₁-C₆alkoxy-C₁-C₆alkyl, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂, C₁-C₆alkylcarbonylamino, C₁-C₆haloalkylcarbonylamino, C₁-C₆alkoxycarbonylamino and C₁-C₆alkylaminocarbonylamino;

R₉ is H, C₁-C₁₂alkyl unsubstituted or substituted by from one to five identical or different substituents selected from halogen, CN, NO₂, OH, C₁-C₆alkoxy, C₁-C₆alkylthio, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ wherein the two alkyl groups are independent of one another and C₁-C₆alkylcarbonylamino; C₃-C₈cycloalkyl, C₁-C₆alkylcarbonyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₂-C₆alkynyl, C₂-C₆haloalkynyl, aryl, heterocycl or benzyl, wherein the aryl, heterocycl and benzyl radicals are unsubstituted or, depending upon the possibilities of substitution, substituted by from one to five substituents selected from the group consisting of halogen, CN, NO₂, OH,

C₁-C₆haloalkyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₃-C₆alkynyl, C₃-C₆cycloalkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₂-C₆alkenylxylo, C₂-C₆haloalkenylxylo, C₃-C₆alkynylxylo, C₃-C₆haloalkynylxylo, C₁-C₆alkylcarbonyl, C₁-C₆haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylthio, C₂-C₆alkenylthio, C₃-C₆alkynylthio, C₁-C₆alkylthio, C₃-C₆haloalkynyl, C₁-C₆haloalkylthio, C₁-C₆alkoxy-C₁-C₆alkyl, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ wherein the two alkyl groups are independent of one another, C₁-C₆alkylcarbonylamino, C₁-C₆haloalkylcarbonylamino, C₁-C₆alkoxycarbonylamino and C₁-C₆alkylaminocarbonylamino;

R₁₀ is H, C₁-C₁₂alkyl unsubstituted or substituted by from one to five identical or different substituents selected from halogen, CN, NO₂, OH, C₁-C₆alkoxy, C₁-C₆alkylthio, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ and C₁-C₆alkylcarbonylamino; C₃-C₆cycloalkyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₂-C₆alkynyl, aryl, heterocycl or benzyl, wherein the aryl, heterocycl and benzyl radicals are unsubstituted or, depending upon the possibilities of substitution, substituted by from one to five identical or different substituents selected from the group consisting of halogen, CN, NO₂, OH, SH, C₁-C₆alkyl, C₁-C₆haloalkyl, C₃-C₆cycloalkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₁-C₆alkylcarbonyl, C₁-C₆haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylcarbonyl-C₁-C₆alkyl, C₁-C₆alkoxy-C₁-C₆alkyl, NH₂, NH(C₁-C₆alkyl), N(C₁-C₆alkyl)₂ wherein the two alkyl groups are independent of one another, C₁-C₆alkylcarbonylamino, C₁-C₆haloalkylcarbonylamino, C₁-C₆alkoxycarbonylamino and C₁-C₆alkylaminocarbonylamino;

R₁₁ and R₁₂ are each independently of the other H, C₁-C₆alkyl, C₁-C₃haloalkyl, C₁-C₆alkylcarbonyl, C₁-C₃haloalkylcarbonyl, C₁-C₆alkoxycarbonyl, C₃-C₆cycloalkyl, C₃-C₆cycloalkyl-C₁-C₆alkyl or C₃-C₆cycloalkylcarbonyl;

R₁₃ is H, C₁-C₆alkyl, C₂-C₆alkenyl, C₃-C₆alkynyl or C₁-C₆haloalkyl;

R₁₄ is H, C₁-C₆alkyl, C₂-C₆alkenyl, C₃-C₆alkynyl or C₁-C₆haloalkyl;

R₁₅ is H, C₁-C₆alkyl, C₂-C₆alkenyl, C₃-C₆alkynyl or C₁-C₆haloalkyl;

k is 0, 1, 2, 3 or 4;

m is 1 or 2; and

q is 0, 1 or 2;

or, where applicable, a possible E/Z isomer, E/Z isomeric mixture and/or tautomer thereof, in each case in free form or in salt form.

2. (Currently amended) The A compound according to claim 1 wherein M is NOR₆.

3. (Currently amended) ~~The A~~ compound according to claim 1 wherein M is O.
4. (Currently amended) ~~The A~~ compound according to claim 1 in free form.
5. (Previously presented) A compound according to claim 1 wherein X₁ and X₂ are chlorine or bromine.
6. (Previously presented) A compound according to claim 1 wherein D is CH.
7. (Previously presented) A compound according to claim 1 wherein A₃ is straight-chain alkylene bridge.
8. (Canceled)
9. (Canceled)
10. (Previously presented) A pesticidal composition which comprises as active ingredient at least one compound defined in claim 1, in free form or in agrochemically acceptable salt form, and at least one adjuvant.
11. (Original) A method of controlling pests which comprises applying a pesticidal composition as defined in claim 10 to the pests or to the locus thereof.